

Modulus Deficit in Thin Film Copper Electrodeposits: Experiments and Modeling

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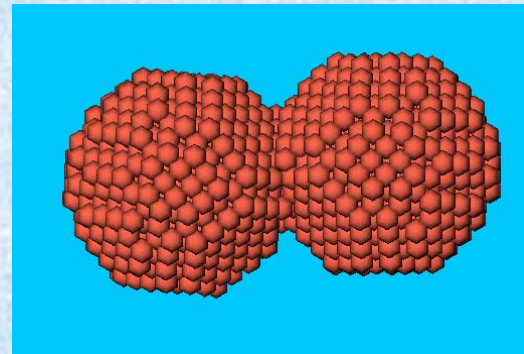
Outline, 1/2

I will describe numerical experiments, based on SEM observations, that may point toward a mechanism for this modulus deficit.

Introduction -- Experiments

SEM observations

Numerical models



Results

Discussion -- What's the mechanism?

Conclusions – Trends of numerical results are in the right direction, and appear large enough to explain the effect – mechanism appears to be “load shedding”

Outline, 2/2

Acknowledgments:

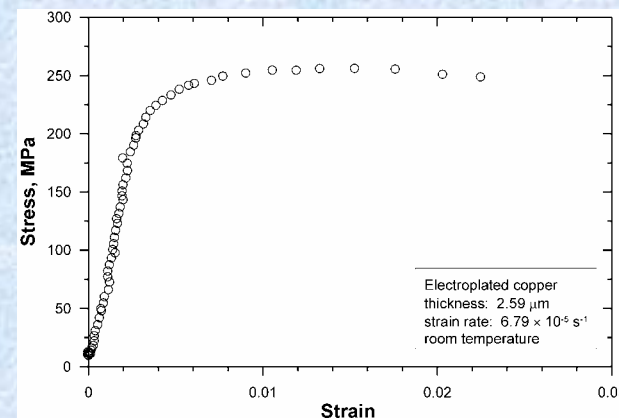
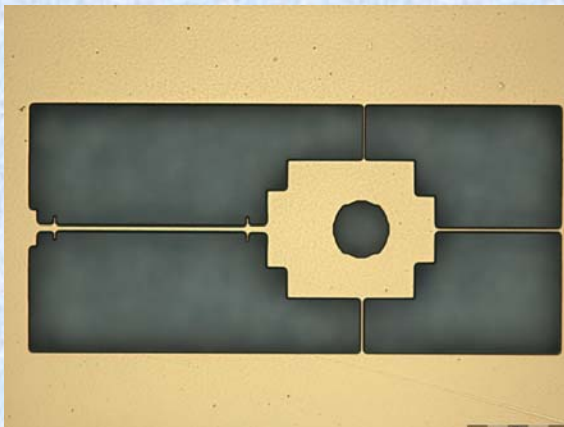
- NIST for funding, including Office of Microelectronics Programs
- Colleagues Roy Geiss for SEM, and Yi-Wen Cheng and Joyce Wright for mechanical testing
- Prof. Jon Rifkin of U. of Connecticut for MD code XMD

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Introduction, 1/2

Microtensile tests of metal thin films (0.5 to 10 μm thick, physical vapor deposited, sputtered, or electrodeposited) generally produce low values of Young's modulus (**10 s of percent**), relative to the bulk polycrystalline average of reference values. (ref Huang and Spaepen, *Acta Met.* **48**, 2000, 3261-3269, my previous results, etc.)

- Microtensile tests done by the same techniques, for some other materials, give the “expected” value: epitaxial Si, CVD Poly Si

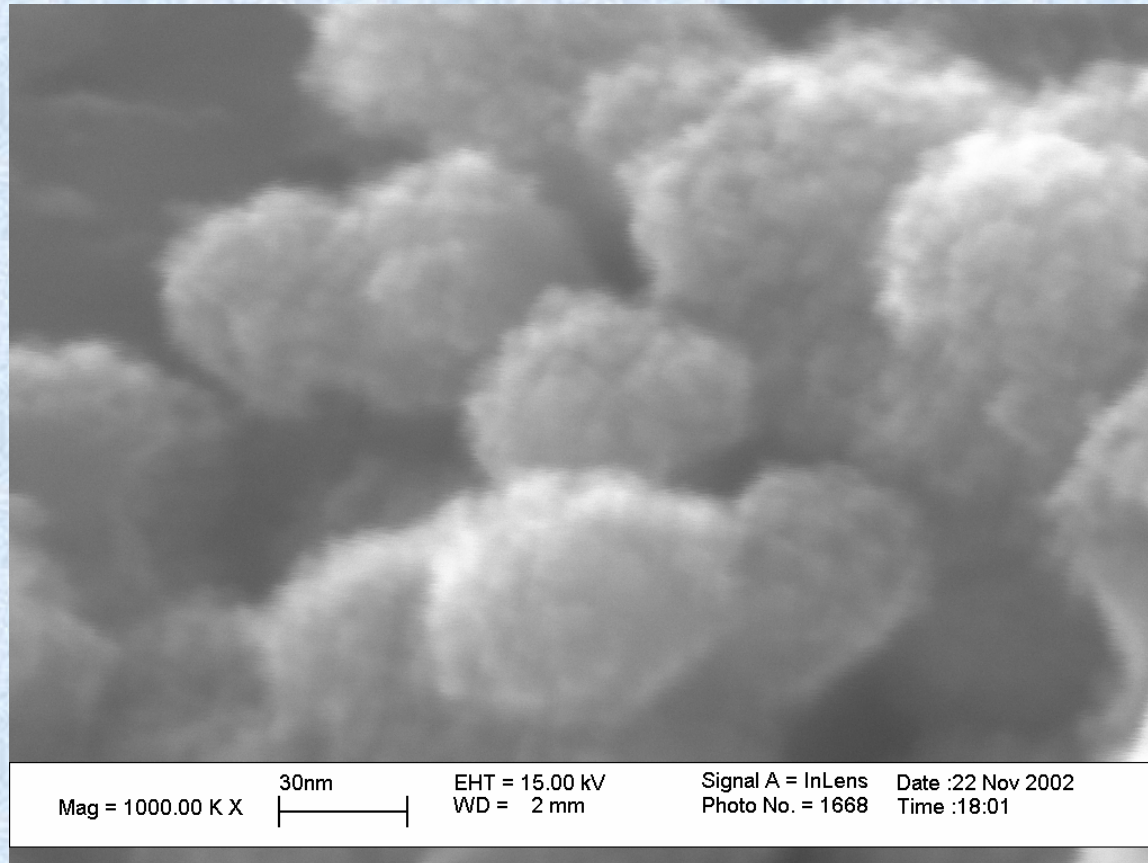


Introduction, 2/2

- A variety of mechanisms for the modulus deficit are suspected, under the general heading of anelasticity.
- Important to understand because:
 - the modulus value is basic to mechanical design;
 - getting the modulus right is basic to mechanical testing.
- Measurements of the elastic properties of thin films have an “interesting” history---these are difficult measurements.

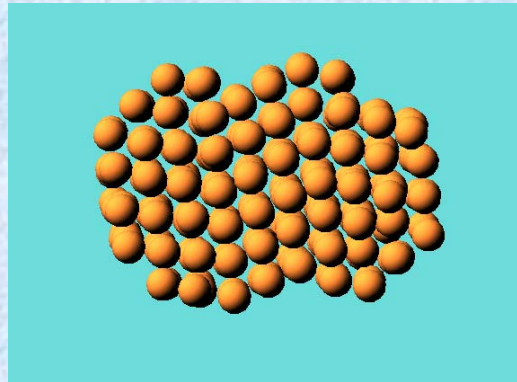
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High-resolution SEM of electrodeposited copper



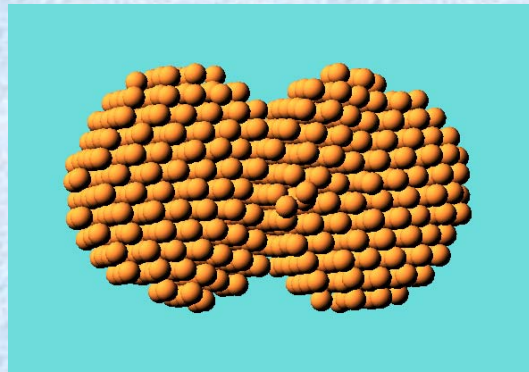
Micrograph at 1 M X clearly shows loosely aggregated spheroids, diameter approximately 30-50 nm, with possible substructure

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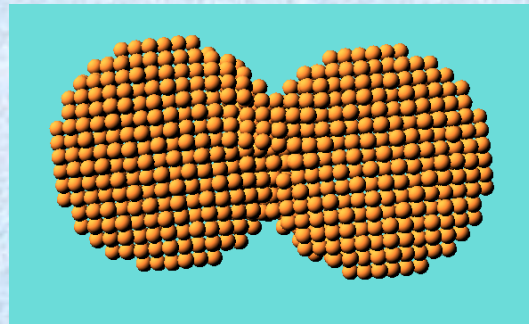


Atoms per initial
sphere

140

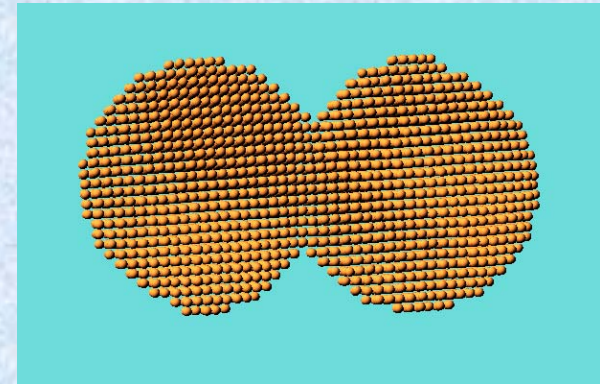


456

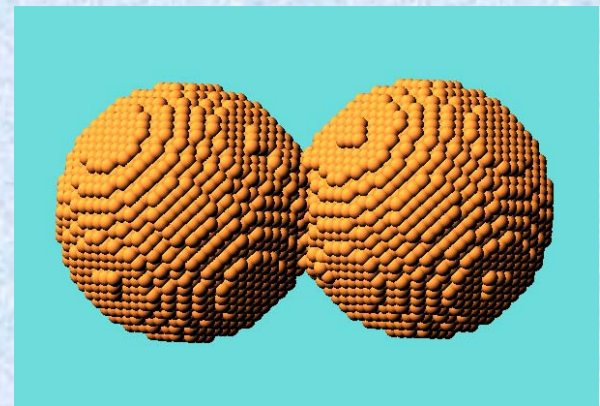


1088

3604



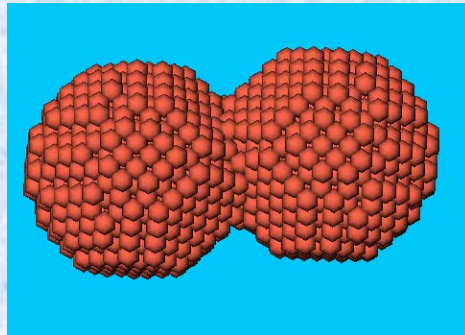
8628



Size effect in metal
sphere aggregation, 0 K

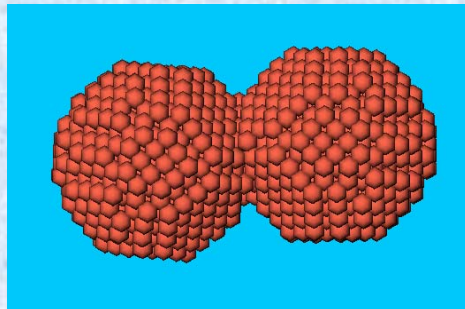
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Results after 9×10^{-12} s (9 picoseconds):

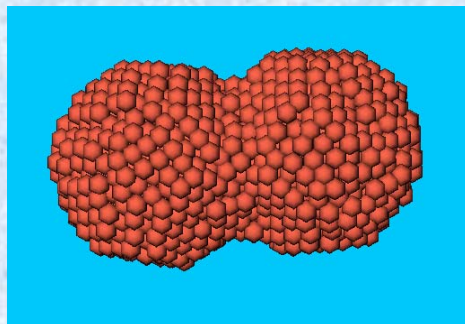


Temp., K:

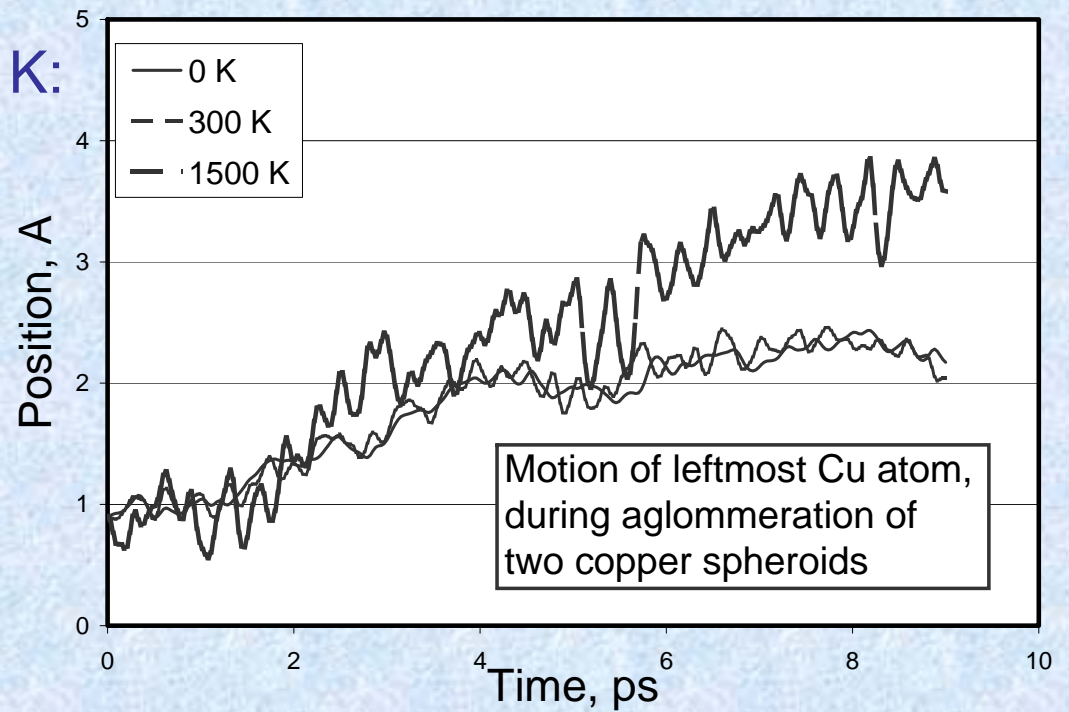
0



300



1500



Result: Spheroids can aggregate at room temperature without electrical forces or solvent effects

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Numerical models: Chains of spheres

3 types:

$\langle 100 \rangle$

$\langle 110 \rangle$

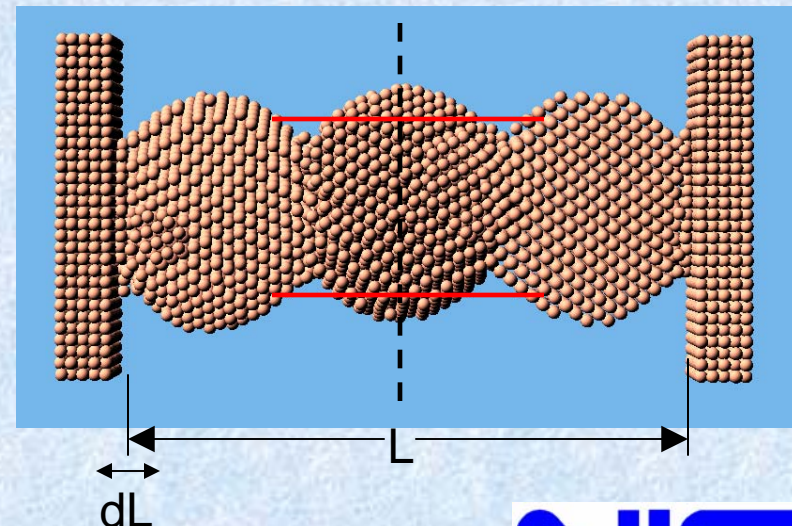
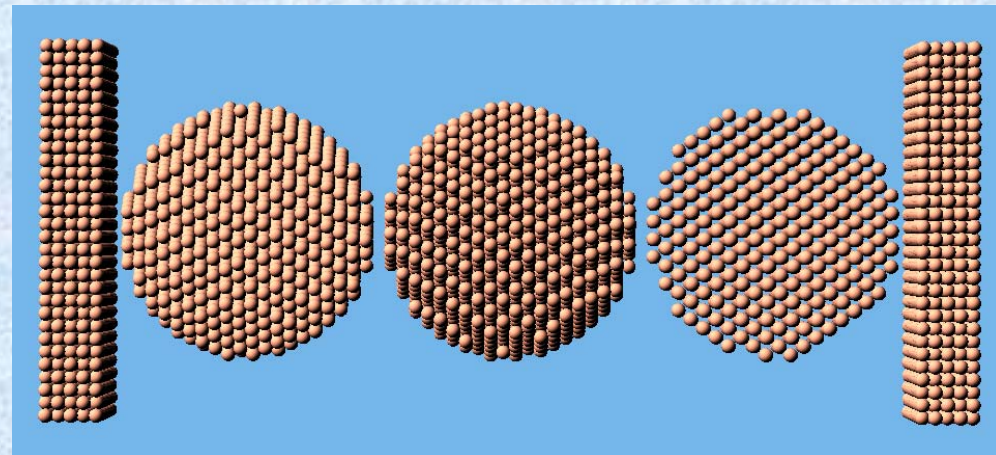
$\langle 111 \rangle$

Along tensile axis

Various twist angles

Around 2000 atoms per sphere, at 300 K

Many-body EAM potential
after Mei and Fernando
(*Phys Rev B* **43**, 1991,
4653-4658)



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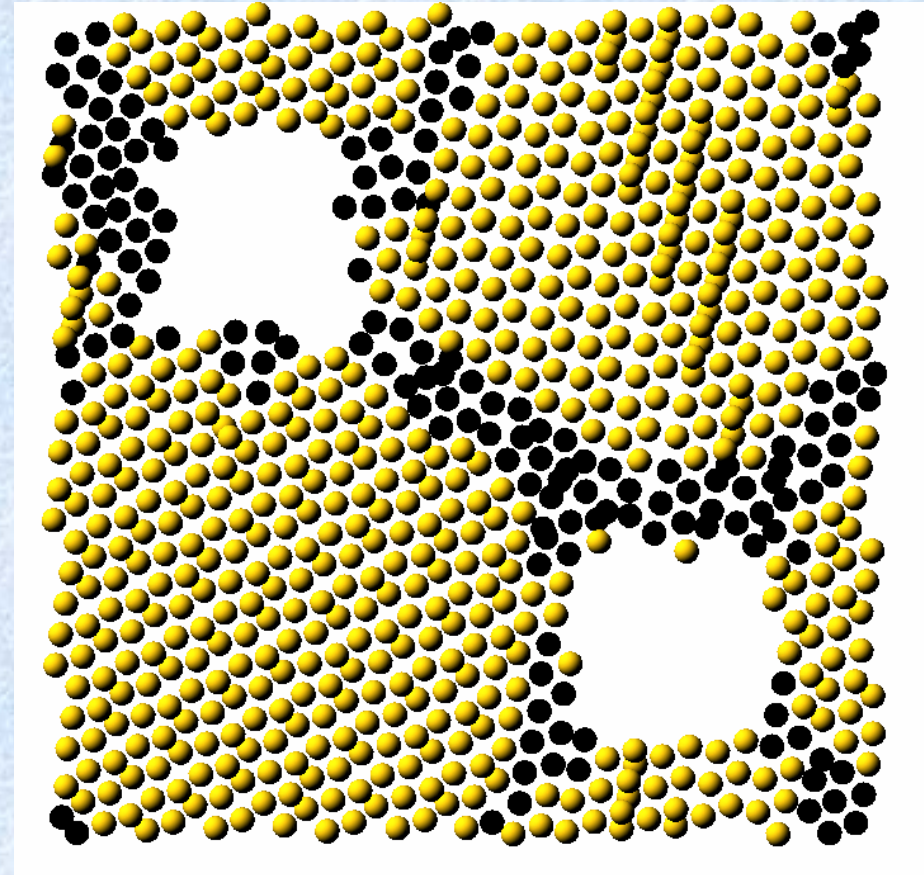
Periodic Boundary
Condition (PBC) model

FCC array of spherical
clusters of atoms

Self-assembled at 300 K

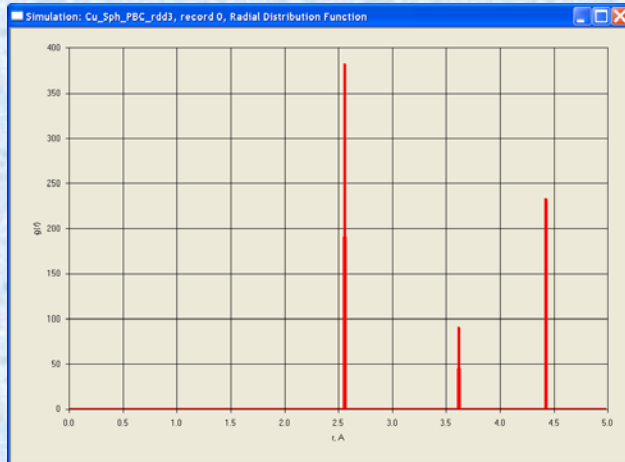
Colored: High FCC
symmetry

Black: Low FCC symmetry



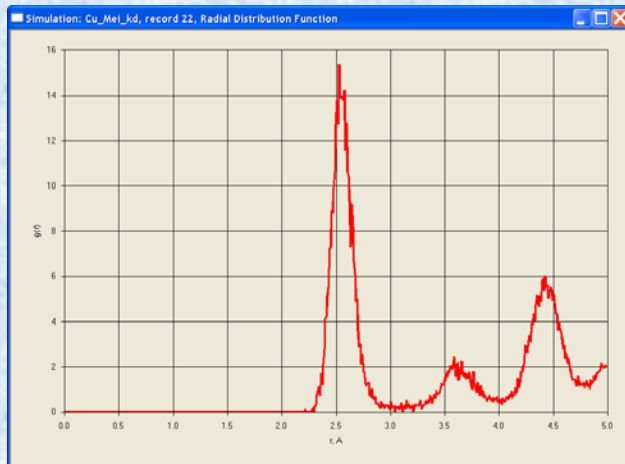
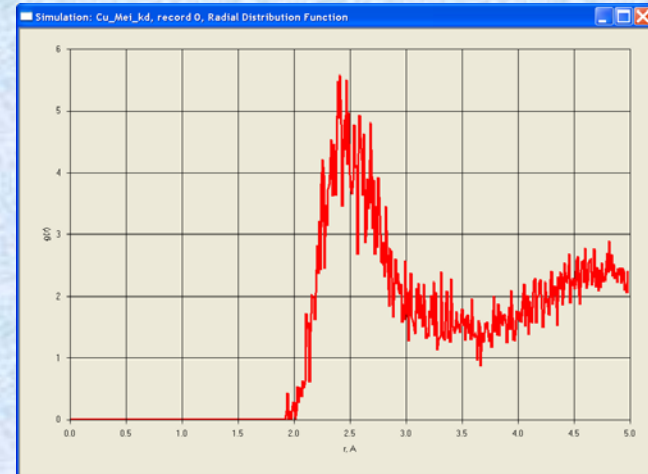
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Radial distribution functions (RDF) for various models



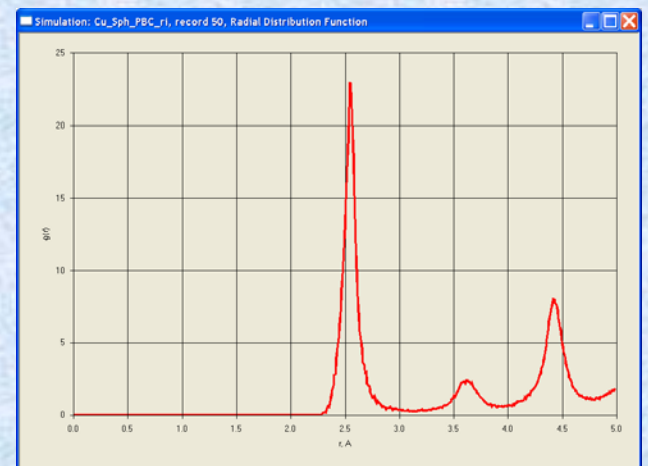
Perfect

Melt temp.
+ 1000 K

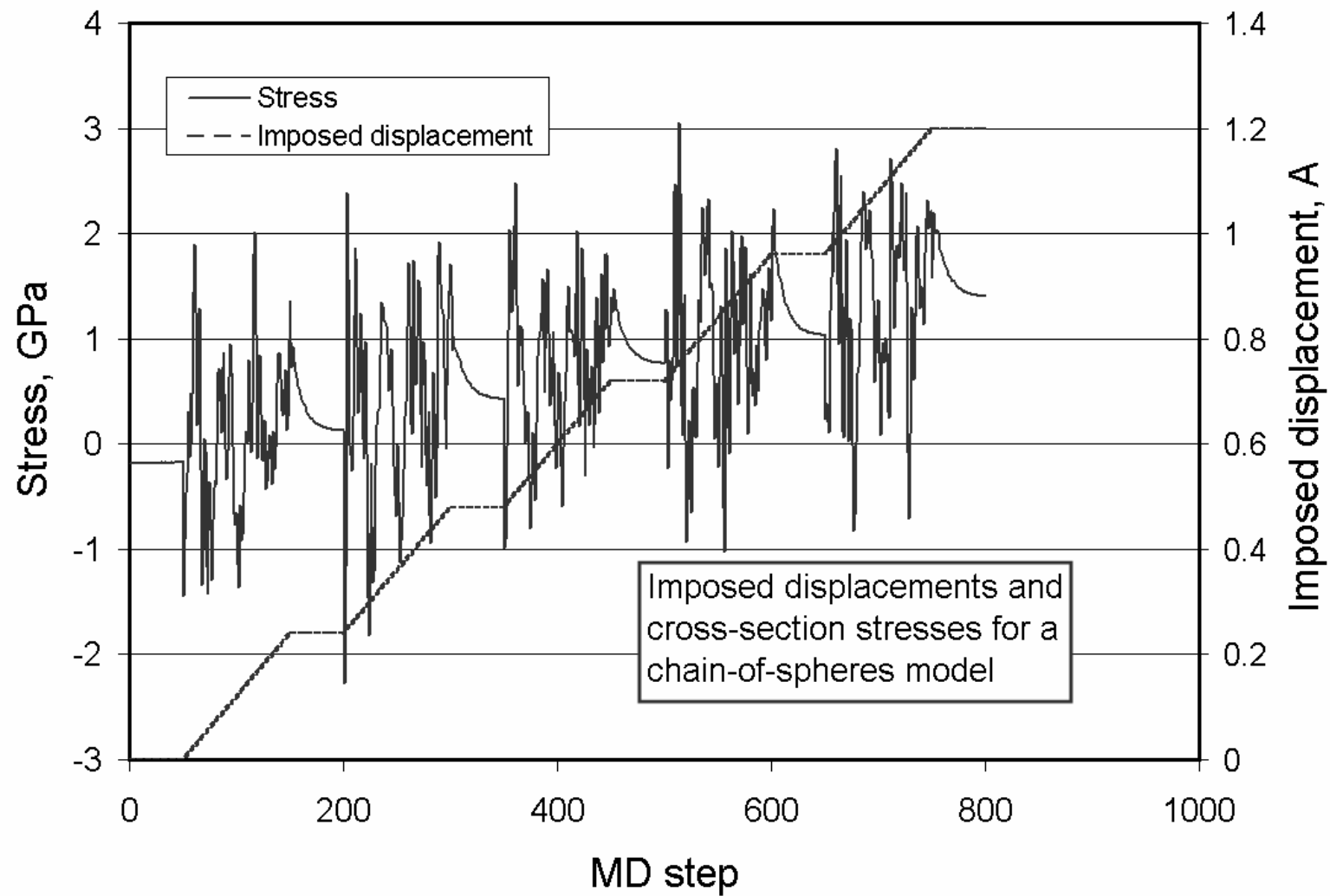


Solidified
block

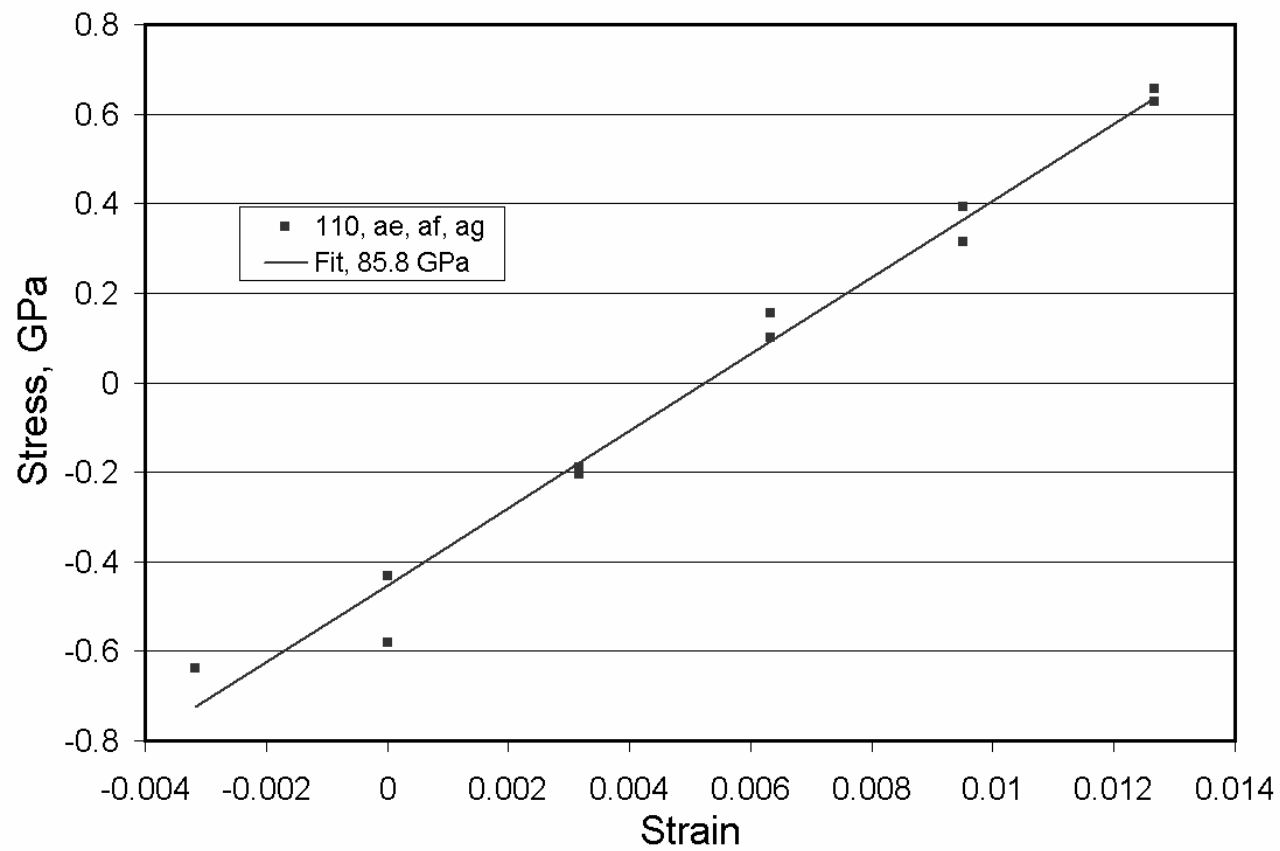
Chain of
spheres



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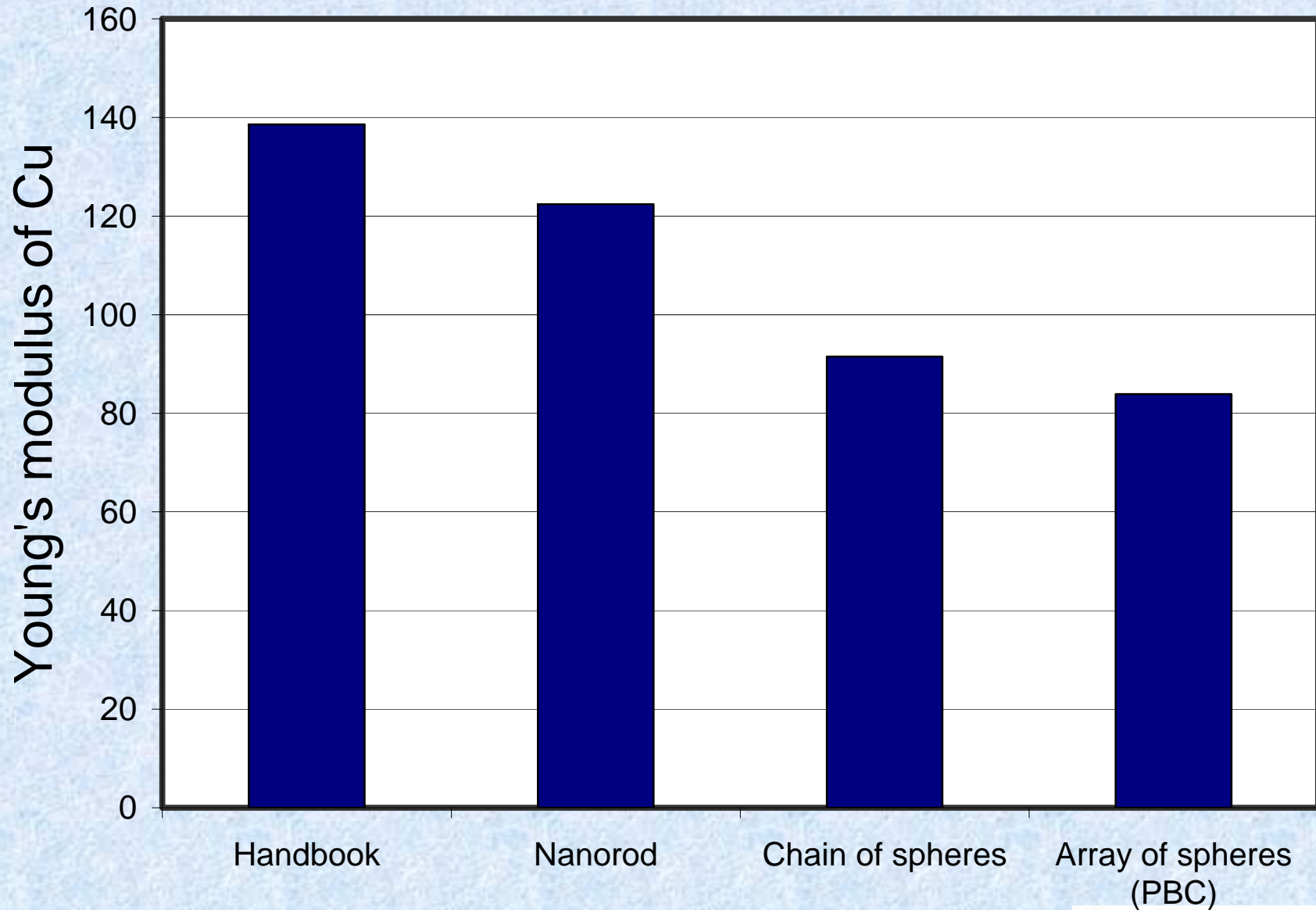
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Stress = F/A , av

Strain = dL/L

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